# ANALYTICAL RESULTS OF SURFACE WATER SAMPLES COLLECTED FROM RACCOON CREEK February 3, 1999 Sampling Event

### Prepared for

### LYONDELL CHEMICAL COMPANY/BEAZER EAST INC.

### Prepared by:

Applied Hydrology Associates, Inc. Pittsburgh, PA Denver, Colorado

February 23, 1999



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4

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### LYONDELL CHEMICAL COMPANY Monaca, Pennsylvania

### TABLE OF CONTENTS

Section	l Pag	<u>je</u>
1.0	INTRODUCTION	
2.0	SAMPLING	
3.0	RESULTS	2
<u>Tables</u>		
1 2	Summary of Analytical Results for Samples Collected from Raccoon Creek	
Figure:	<u>s</u>	
1 2	Transect Location Map Surface Water Benzene Concentrations at Transect E	<b>4 5</b>
Appen	<u>dices</u>	
A B	Groundwater Elevations, East and West Sides of Raccoon Creek Data Validation Report	

### LYONDELL CHEMICAL COMPANY Monaca, Pennsylvania

#### 1.0 INTRODUCTION

This report presents the results of surface water samples collected from Raccoon Creek at the Lyondell Chemical Company (LCC) / Beazer East Inc. (BEI) Monaca, PA site during the February 3, 1999 quarterly monitoring event. The samples were collected in compliance with Appendix D of the 1997 Consent Order and Agreement (1997 CO&A) between ARCO Chemical Company<sup>1</sup>, BEI and the Pennsylvania Department of Environmental Protection (PADEP) dated October 20, 1997.

#### 2.0 SAMPLING

Surface water samples were collected at Transect E as defined in the 1997 CO&A. The locations of Transect E is shown in Figure 1. In addition, water elevations were measured in nearby monitoring wells and the results are presented in Appendix A.

A total of eight surface water samples, including a duplicate were collected from Raccoon Creek on February 3, 1999<sup>2</sup>. These samples were collected at the same three locations along Transect E as in previous sampling events. The locations are shown in Figure 2 and are at the center of the stream, and approximately 30 feet from the east and west banks. At the center location, samples were collected at three depths; 6 inches below the surface, 2 inches above the bottom, and midway between the surface and bottom. Samples from the east and west sides of the transect were collected at two depths; 2 inches above the bottom, and midway between the surface and bottom.

During sampling a boat was stationed at Transect E using a rope secured to the east and west shores of Raccoon Creek. The samples were collected by using a peristaltic pump to pump water from the desired depth into three 40-ml vials preserved with hydrochloric acid. Samples were collected from the required depths utilizing tubing secured to a vertical steel rod lowered from the boat until it rested on the bottom of the creek. The rod did not penetrate the sediments on the creek bottom because a 1-foot diameter disc constructed of steel mesh is fastened perpendicular to the bottom of the rod.

Two tubes were used. The bottom of the "deep sample tube" was secured to the probe 2 inches from the bottom of the probe. The bottom of the "mid-depth sample tube" is adjustable and was secured to the probe mid-depth at each location. Care was taken not to disturb the sediments at the sampling location and the pumped water was closely monitored to ensure sediment was not included in the sample. One gallon of water was pumped through the tubing before each sample is obtained in order to purge the tubing.

ARCO Chemical Company is now Lyondell Chemical Company

<sup>&</sup>lt;sup>2</sup> A field blank was also obtained after pumping one gallon of distilled water through the tubing used for sampling.

Monaca, Pennsylvania

The samples were uniquely numbered as follows to identify the location, depth and date of sampling:

RC-EC-00-0299

Where:

RC indicates Raccoon Creek
 EC indicates Transect E and location (C = center, L = left bank, R = right bank [facing downstream])
 indicates sample depth in feet and tenths of a foot (0.0 feet)
 indicates the month and year collected (February 1999)

Samples were logged onto a chain of custody form (included in of the data validation report in Appendix B) and stored on ice until receipt by Reliance Laboratories Inc. in Edison, NJ. Reliance analyzed the samples using USEPA Method 524.2 for BTEXS.

#### 3.0 RESULTS

The analytical results are presented in Table 1. Benzene was detected in six of the eight samples and concentrations in samples where benzene was detected ranged from  $0.37~\mu g/L$  in Sample RC-EL-18-0299 to  $0.69~\mu g/L$  in sample RC-EC-64-0299. Sampling locations and depths are shown on Figure 2, along with the concentration of benzene at each location. Water levels in wells near Raccoon Creek are presented in Appendix A.

Table 1
Summary of Analytical Results for Samples Collected from Raccoon Creek

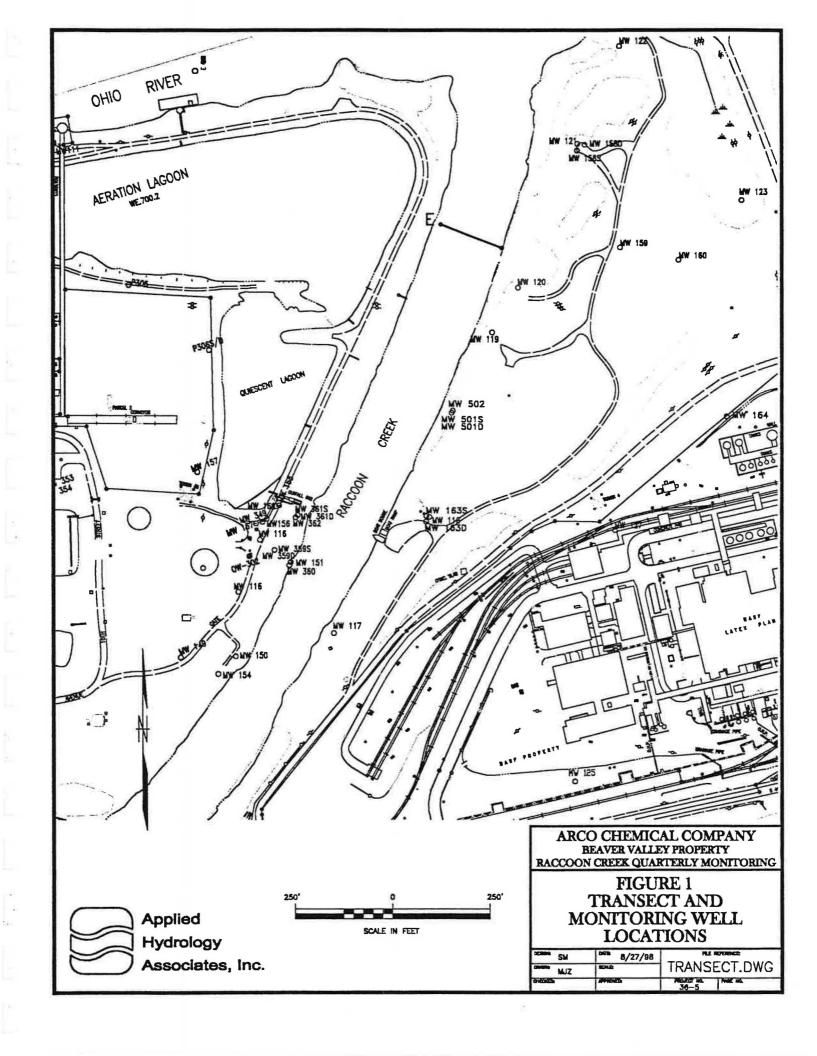
Sample Name	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
RC-EL-18-0299	0.37	0.78	0.27	1.17	< 0.58
RC-EL-36-0299	0.49	0.82	0.32	1.37	< 0.58
RC-EC-00-0299	0.58	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0299A	0.64	< 0.6	< 0.22	0.61	< 0.58
RC-EC-34-0299	0.64	0.61	< 0.22	0.75	< 0.58
RC-EC-64-0299	0.69	< 0.6	< 0.22	0.55	< 0.58
RC-ER-12-0299	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-25-0299	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58

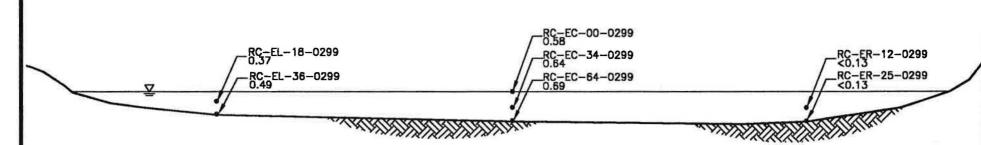
The analytical data were validated upon receipt and found to be acceptable. A Data Validation Report which includes the Certificate of Analysis is provided as Appendix B. Table 2 presents the historical concentration of benzene in Raccoon Creek at Transect E during all monitoring events to date.

Table 2
Historic Benzene Concentrations at Transect E
(ug/L)

Sampling Location	Sampling Depth	7/23/97	10/28/97	2/25/98	5/21/98	7/29/98	10/27/98	2/3/99
30 Feet off West Bank	Mid-depth	0.28	<0.13	<0.13	0.70	<0.13	1.57(1)	0.37
30 Feet off West Bank	Deep	0.81	<0.13	<0.13	0.70	<0.13	0.61 <sup>(1)</sup>	0.49
Center of Creek	Shallow	0.24	<0.13	0.38	0.70	<0.13	<0.13	0.61(1)
Center of Creek	Mid- Depth	0.18	<0.13	0.49	0.64	<0.13	0.2	0.64
Center of Creek	Deep	0.46	<0.13	0.30	0.60	<0.13	<0.13	0.69
30 Feet off East Bank	Mid-depth	0.16	<0.13	<0.13	<0.13	0.13	0.52	< 0.13
30 Feet off East Bank	Deep	<0.13	<0.13	0.14	0.22	0.22	<0.13	< 0.13

<sup>(1)</sup> Results shown are the average of the blind duplicate samples.





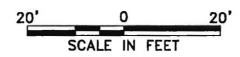
## CREEK SECTION LOOKING DOWNSTREAM

### **LEGEND**

SURFACE WATER SAMPLE LOCATION
 ALL CONCENTRATIONS IN ug/L



Applied Hydrology Associates, Inc.



LYONDELL CHEMICAL COMPANY
BEAVER VALLEY PROPERTY
RACCOON CREEK QUARTERLY MONITORING

FIGURE 2

SURFACE WATER
BENZENE CONCENTRATIONS
AT TRANSECT 'E'

**FEBRUARY 3, 1999** 

SM SM	8/27/98	BENZENE.dwg			
MJZ	NOT TO SCALE				
D-Marie	AMORE	36-5 ME IS.			

### Appendix A

# Groundwater Elevations, East and West Sides of Raccoon Creek

# GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK February 3, 1999

Well Number	Top of Casing (TOC) (ft. amsl)	Depth to SPL from TOC (2) (ft. amsl)	Depth to Water from TOC (2) (ft. amsl)	Calculated Water Level Elevation (1) (ft. amsl)	Calculated SPL Thickness (3) (ft. amsl)	Comments
		Mo	onitoring We	ells Screened in	Silty Clay U	Jnit
MW 260	(05.04	) ID	2.02	OTH AREA	NY/A	
MW - 360	685.84	ND	2.03	683.81	N/A	
MW - 170	706.70	ND	22.02	684.68	N/A	
MW - 362	689,43	ND	5.59	683.84 CCOON CREEK A	N/A	
B/TW 110	(00.20	) ID			N/A	
MW- 118	690.39	ND	6.88	683.51		
MW - 502	701.86	ND	18.36	683.50	N/A	
MW - 119	705.59	ND	22.08	683.51	N/A	
MW - 120	709.42	ND	25.85	683.57	N/A	
MW - 121	713.90	ND	30.33	683.57	N/A	
MW - 152	696.35	ND	12.82	683,53	N/A	
		Monitorin	g Wells Scre	eened in Upper	Sand and G	Gravel Unit
				OTH AREA		
MW - 344	709.42	ND	25.41	684.01	N/A	
MW - 359S	692.93	ND	9.24	683.69	N/A	
MW - 361S	689.40	ND	5.80	683,60	N/A	
MW - 169	707.93	ND	24.22	683.71	N/A	
MW - 167	711.06	ND	27.35	683.71	N/A	Top of casing changed from 707.36 to 711.06 on 11/98 to accommodate respiration monitoring well head. Monitoring well stick up is 3.70 above orig TOC
				CCOON CREEK A		
MW - 163S	690.87	ND	7.32	683,55	N/A	
MW - 501S	701.30	ND	18.05	683.25	N/A	
MW - 162S	706.05	ND	22.54	683.51	N/A	
MW - 159	708.99	ND	25.43	683.56	N/A	
MW - 160	701.00	ND	17.48	683.52	N/A	
MW - 158S	713.60	ND	30.04	683.56	N/A	
MW - 122	692.78	ND	9.23	683.55	N/A	
Note: See figure						
(1) Calculated	values, based or	Elevation of To	OC minus Depth t	o Water from TOC.		
(2) Measured f	rom top of casin	g using the MM	A Interface Probe	. ND means no SPL	was detected.	
(3) Calculated	values, based or	Depth to Water	from TOC minu	s Depth to SPL from	TOC. N/A means	not applicable, no SPL was detected.

## GROUNDWATER LEVELS ON THE EAST AND WEST SIDES OF RACCOON CREEK February 3, 1999

Well Number	Top of Casing (TOC) (ft. amsl)	Depth to SPL from TOC (2) (ft. amsl)	Depth to Water from TOC (2) (ft. amsl)	Calculated Water Level Elevation (1) (ft. amsl)	Calculated SPL Thickness (3) (ft. amsl)	Comments	
		Monitorin	g Wells Scre		Sand and Grave	el Unit	
				OTH AREA			
MW 345	708.91	ND	25.37	683.54	N/A		
MW 361D	689.35	ND	5.75	683,60	N/A		
MW 359D	692.80	ND	9.23	683.57	N/A		
				CCOON CREEK A			
MW 163D	689.62	ND	6.03	683.59	N/A		
MW 501D	701.44	ND	18.00	683.44	N/A		
MW 166D	703.95	ND	20.51	683.44	N/A		
MW 158D	712.04	ND	48.70	663.34	N/A		
		Wat	er Levels in	Raccoon Creel	k and Ohio Rive	r	
			RACCOON	CREEK AREA STA	FF GAUGE		
Time of Observation	Staff Gauge Elevation (a) (ft. amsl)	Staff Gauge Reading	Calculated Water Level Elevation (ft. amsl)	Comments			
8:13	685.00	1.31	683.31				
9:16	685.00	1.40	683,40			10.00	
7.10	005,00	1		RIVER, STAFF GA	AUGE		
9:04	685,96	3,40	683,36				
10:00	685,96	3.50	683.46				
					4		
Note: See figur					<u> </u>		
				to Water from TOC.			
				e. ND means no SPL s Depth to SPL from		pplicable, no SPL was detected.	
(4) Elevation 6	85.00 is equival	ent to 3.00 mark	on staff gauge at	Raccoon Creek			
(5) Elevation 6	85,96 is equival	ent to 6.00 mark	on staff gauge at	Ohio River			

### Appendix B

**Data Validation Report** 

AHA Filename: BVRC798.doc



1200 South Parker Road, Suite 100

Denver, CO 80231

Tel: (303) 873-0164

Fax: (303) 873-6110

### MEMORANDUM

TO:

**Files** 

FROM:

Skip Meier, Applied Hydrology Associates

DATE:

February 22, 1999

**SUBJECT:** 

Data Validation Results, Lyondell Chemical Company Beaver Valley Property

Data validation was performed on the volatile organic analytical data from nine surface water samples obtained from Raccoon Creek on February 3, 1999. The validation was performed in accordance with the "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review". Reliance Laboratories Inc. performed the analysis using EPA Method 524.2. The samples reviewed included:

Field Sample ID	Lab Sample ID
RC-EL-18-0299	R6171.5
RC-EL-36-0299	R6171.4
RC-EC-00-0299	R6171.6
RC-EC-00-0299A	R6171.7
RC-EC-34-0299	R6171.8
RC-EC-64-0299	R6171.3
RC-ER-12-0299	R6171.1
RC-ER-25-0299	R6171.2
Field Blank	R6171.9

Items reviewed and actions taken were as follows:

#### √ Method:

The nine samples were analyzed for BTEXS by method USEPA 524.2 on February 4, 1999.

#### ✓ Holding Time:

All Samples were analyzed within the 14-day holding time.

#### √ Blanks

No target compounds were detected in the associated method blank.

### **√** System Monitoring Compounds:

The "USEPA Contract Laboratory Program National Functional Guidelines for Organic Data Review" indicate that "Recoveries for system monitoring compounds in volatile samples and blanks must be within the limits specified in the Method." However, Method 524.2 does not specify a required recovery. Nevertheless, 4-bromofluorobenzene and 1,2-dichlorobenzene-d4 surrogate recoveries were within 95-101 percent and this is acceptable.

### √ Internal Standards:

All fluorobenzene internal standards were within the established criteria for area internal standard and retention time.

### **V** GC/MS Instrument Performance Check:

All bromofluorobenzene (BFB) tunes met the ion abundance criteria. Analysis of the instrument performance check solution was performed at the beginning of each 12-hr period during which the samples were analyzed.

#### **√** Initial Calibrations:

The initial calibration performed on February 4, 1999 for Instrument HP5971A met the 30 percent relative standard deviation (RSD) and 0.05 minimum relative response factor criteria for all compounds.

### **√** Continuing Calibrations:

Continuing calibration was run and compared to the correct initial calibration. All continuing calibrations met the 25 percent difference and minimum relative response factor criteria for all compounds.

### √ Matrix Spike/Duplicate:

The matrix spike/duplicate results for recovery and RPD were within the Quality Control limits.

### **√** Target Compound Indentification/Quantitation:

No problems were identified with compound identification or quantities.

### √ Field Duplicate:

A field duplicate was collected during this sampling event. The duplicate sample was denoted by an "A" at the end of the sample name. The pair is RC-EC-00-0299 and duplicate RC-EC-00-0299A. Table 1 below summarizes the RPD for the sample/duplicate pair.

Table 1: Relative Percent Difference (RPD)

	Sample Name	Benzene (ppb)	RPD (%)	Toluene (ppb)	RPD (%)	Ethyl- Benzene (ppb)	RPD (%)	Xylene (ppb)	RPD (%)	Styrene (ppb)	RPD (%)
Ì	RC-EC-00-0299	0.58	8.3	ND	NA	ND	NA	ND	NA	ND	NA
	RC-EC-00-0299A	0.64		ND	NA	ND	NA	0.61	NA	ND	NA

ND = Non Detect

NA = Not Applicable

#### √ Summary:

The overall quality of the data was good. There was good agreement between duplicate sample pair RC-EC-00-0299 and RC-EC-00-0299A (See Table 1). The concentration of benzene in the field blank was 0.63 ug/L, but this was attributed to lab error since one gallon of distilled water was pumped through the sampling tubing before taking the field blank. No trip blank was included with the samples sent to the lab. This oversight will be remedied in subsequent sampling events.

### RELIANCE LABORATORIES, INC.



175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

### **ANALYTICAL REPORT**

For Arco Chemical Co. Pittsburgh, PA 15219

Project: Arco / Monaca

### RELIANCE LABORATORIES, INC.



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### **ANALYTICAL DATA REPORT**

for

Arco Chemical Co. Pittsburgh, PA 15219 Project: Arco / Monaca

Date Received: 2/4/99

Sample ID	Lab ID #
Sample ib	
RC-ER-12-0299	R-6171.1
	R-6171.2
RC-ER-25-0299	R-6171.3
RC-EC-64-0299	R-6171.4
RC-EL-36-0299	R-6171.5
RC-EL-18-0299	1
RC-EC-00-0299	R-6171.6
RC-EC-00-0299A	R-6171.7
RC-EC-34-0299	R-6171.8
	R-6171.9
Field Blank	

These samples have been analyzed by EPA method 524.2 for a selected compound list. The results are not designed for use for drinking water purposes.

G. P. Kirpalani Manager

GPK/vb

### RELIANCE LABORATORIES INC.



### 175 MAY STREET, EDISON, NJ 08837 PH (732) 738-5454 FAX (732) 738-5841 EMAIL: 74201.3501@COMPUSERVE.COM

### TABLE OF CONTENTS

Laboratory Chronicle	page	1
Non-Conformance Summary	page	2
Methodology Summary	page	3
Analytical Results	page	5
QA/QC Data	page	6
Laboratory Certificate	page 3	33
Chain of Custody	page 3	34

### RELIANCE LABORATORIES, INC.



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### LABORATORY CHRONICLE

Customer Name Arco Chemic Date Received: 2/4/99 Date Sampled: 2/3/99 Sample ID: As per chain of cu		
Organic Extraction:		
1 Acids 2 Base / Neutrals		
3 Pesticides/PCB's		
Analysis:		
1 Volatiles	2/4/99	
A D 01 1 1		
A Destinides (DODIs		
5 TPHC		
Inorganics:		
1 Metals		
3 Phenols		
Other Analysis:		
Supervisor Review & Approval	Geknell	

RELIANCE LABORATORIES INC.



3090 WOODBRIDGE AVENUE, EDISON NJ 08837 PH (908) 738-5454 FAX (908) 738-5841

### **NON-CONFORMANCE SUMMARY**

Reliance Labs received 9 water sample for BTEXS (EPA 524.2) from Arco Chemical on 4 February 1999. Samples consisted of 9 vials.

Matrix spike recovery analysis was performed on samples and results are attached.

All analyses were performed within the required holding time.

### STANDARD OPERATING PROCEDURE METHOD 524.2

1. Scope

This is the general method for the procedure used to identify purgeable volatile organics in portable water. The sample is purged with ultra high purity helium and concentrated into a trap. The volatiles are then thermally desorbed onto a megabore column and identified using a mass spectrometer detector.

- 2. Equipment and Apparatus
- A. Sample containers- 40ml screw caps amber vials.
- B. Purge and Trap System.
  - 1. 25cm VOCARB 3000 trap.
- C. Glassware
  - 1. 20 ml fritted purging vessels.
  - 2. 25 ml teflon sealed syringe with lever lock assembly.
  - 3. 10 μL syringes.
- D. Gas Chromographic / Mass Spectrometer.
  - 1. Column type J&W

75 m, 0.53 mm ID, DB624 3 microns

- E. Apparatus Conditions
  - 1. Tekmar (purge and trap)
    - a. Purge time

2 min.

b. Desorb time and temp.

250° for 2 min.

c. Bake time and temp. :

260° for 12 min.

d. Flow rate

15 cc/min.

- 2. GC Conditions
  - a. Column flow

15 cc/min.

b. Initial temp.

35° C

c. Ramping Rate

6° C/min.

d. Final temp.

200° C

e. Run time

47.25 min.

f. Initial time

6 min.

- 3. Stock Standards
- A Internal Standard
  - 1. Flourobenzene
- B. Surrogates
  - 1. 1.2-dichlorobenzene-d4
  - 2. 4-bromoflurobenzene
- C. Prepare standard solutions for all target compounds and surrogates at 20 ppm.
- D. Prepare internal standard at 20 ppm in methanol.
  - 1. Prepare all standards and store in teflon sealed 1 ml vials.

### 4. Run Sequence

- A. Tune Instrument
- 1. Inject 1µL of 25 ppm BFB into GC.
  - a. Tune must pass against criteria.
  - b. Tune must be run before any samples, blank or calibrations can be run.
  - c. From time to tune 12 hours are available to run all QC data and samples.

### B. Five Point Calibration Curve

- 1. Purge five (5) concentrations of standard solutions containing all the target analysis at 1 ppb, 2 ppb, 5 ppb, 10 ppb, and 20 ppb.
- 2. The above standard must be run within 12 hours of injecting the BFB tune.
- Created a calibration curve with the above standard runs.
  - a. If the 30% RSD deviation is exceeded the standards must be run again (still within 12 hours)
- 4. Create an identification file from this calibration curve for automated quantification.
- C. If time remains in the 12-hour run period go to step F.
- D. If the 12-hour period has expired, a new tune must be injected and a new sequence must be started.
- E. Once an initial calibration curve is established a continuing calibrations check may be run. A continuing calibration check is required every time the mass spectrometer is tuned.
  - 1. 2 ppb concentration of all target compounds is purged and quanted against current ID file.
  - 2. Check the response factors of this run against the average RF of the calibration file. The RF of the continuing calibration must be within ± 50% D (difference) of the 5 point for all compounds.
  - 3. The area counts of internal standard and surrogates must not be decreased by >30% from the most recent continuing calibration standard nor decrease by >50% from the initial calibration standard.

### F. Daily Blank

- 1. Purge 20 ml of laboratory reagent water (nanopure) with 5 ppb internal standard and 5 ppb each surrogate.
- Run this blank and quant against current ID file.
- 3. If blank does not meet criteria, it must be rerun before analyzing any samples.

### G. Samples

- 1. Fill 25 ml syringe until it overflows with sample. Then adjust the volume to 20 ml exactly.
- 2. Inject 5 µl each 25 ppm internal standard and surrogate standard solution into each sample.
- 3. Run and quant against the current 5 point calibration curves.
- Any sample with target compound over 50 ppb must be rerun at the appropriate dilution.
- 5. Any sample not injected in 12-hour period must be rerun.

### H. Quality Control Sample (QCS)

1. Analyze a QCS from an external source at least quarterly.



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LABORATORY ID NJ DEP NO. 12687 PA DER NO. 68437

### **CERTIFICATE OF ANALYSIS**

**Customer:** 

Arco Chemical

Sample:

**Aqueous Samples** 

Date Sampled:

3 February 1999

Lab ID:

R-6171

Reference:

AHA / Monaca

19 February 1999

Units: μg/L

Sample ID	Benzene	Toluene	Ethylbenzene	Xylene	Styrene
RC-ER-12-0299	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-ER-25-0299	< 0.13	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-64-0299	0.69	< 0.6	< 0.22	0.55	< 0.58
RC-EL-36-0299	0.49	0.82	0.32	1.37	< 0.58
RC-EL-18-0299	0.37	0.78	0.27	1.17	< 0.58
RC-EC-00-0299	0.58	< 0.6	< 0.22	< 0.22	< 0.58
RC-EC-00-0299A	0.64	< 0.6	< 0.22	0.61	< 0.58
RC-EC-34-0299	0.64	0.61	< 0.22	0.75	< 0.58
Field Blank	0.63	< 0.6	< 0.22	< 0.22	< 0.58

G. P. Kirpalani Manager

Data File: c:\hpchem\1\data\v5969.d

Vial: 10 Acq On : 4 Feb 99 4:39 pm Operator: vb

Inst : 5971 - In Multiplr: 1.00

Sample : R-6171.1 Misc : Arco - RC-ER-12-0299

Quant Time: Feb 5 9:11 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units Dev(	Min)
1) Fluorobenzene	12.99	96	1967274	5.00 ug/L 0	.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.08 31.26	95 152	701667 446627	J.	ery .07% .76%

Target Compounds

Qvalue

<sup>(#) =</sup> qualifier out of range (m) = manual integration v5969.d RUN524.M Fri Feb 05 09:18:35 1999

Vial: 10

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : c:\hpchem\1\data\v5969.d

Acq On : 4 Feb 99 4:39 pm

Sample : R-6171.1

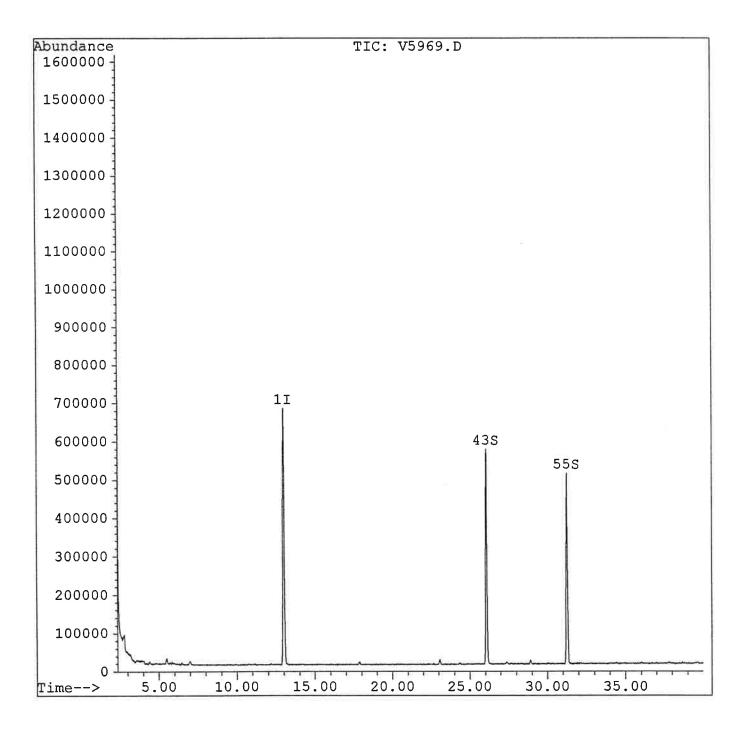
Misc

: Arco - RC-ER-12-0299

Quant Time: Feb 5 9:11 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



Data File : c:\hpchem\1\data\v5970.d

Acq On : 4 Feb 99 5:24 pm Sample : R-6171.2 Misc : Arco - RC-ER-25-0299

Quant Time: Feb 5 9:11 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units De	ev(Min)
1) Fluorobenzene	12.99	96	1877487	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.07 31.25	95 152	661048 447619	%Re 4.84 ug/L 4.98 ug/L	covery 96.81% 99.51%

Target Compounds

Qvalue

Vial: 11

Inst : 5971 - In

Operator: vb

Multiplr: 1.00

<sup>(#) =</sup> qualifier out of range (m) = manual integration v5970.d RUN524.M Fri Feb 05 09:18:43 1999

Vial: 11

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : c:\hpchem\1\data\v5970.d

Acq On : 4 Feb 99 5:24 pm

Sample : R-6171.2

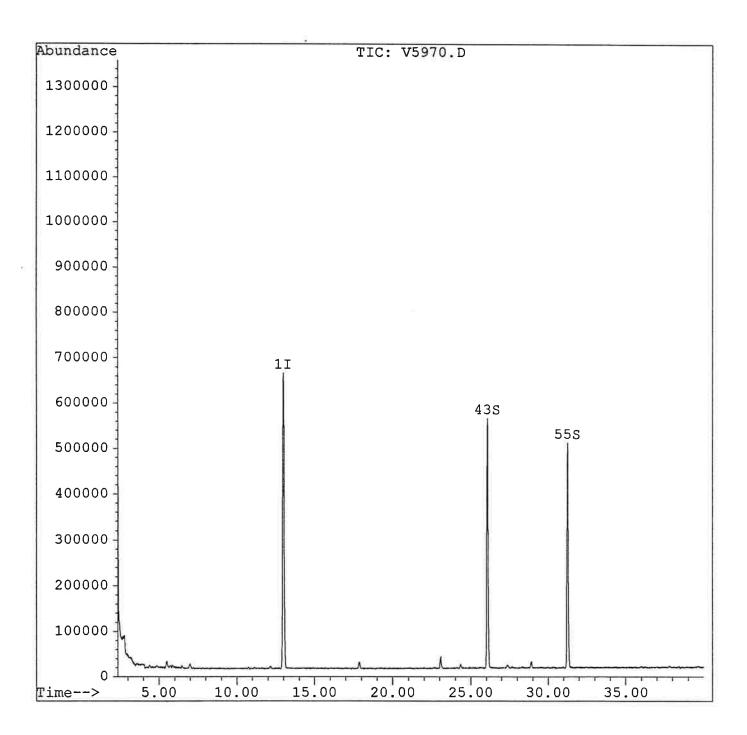
Misc

: Arco - RC-ER-25-0299

Quant Time: Feb 5 9:11 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V5971.D

Vial: 12

Acq On : 4 Feb 99 6:10 pm

Operator: vb

: R-6171.3 Sample

Inst : 5971 - In

Misc : Arco - RC-EL-64-0299

Multiplr: 1.00

Quant Time: Feb 5 9:25 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.99	96	1941054	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.06 31.25	95 152	676366 459698	%F 4.79 ug/L 4.94 ug/L	Recovery 95.81% 98.85%
Target Compounds 19) Benzene 26) Toluene 36) m&p-xylenes 37) o-xylene	12.16 17.84 23.05 24.34	78 91 106 91	254867 161641 71868 79167	0.69 ug/L 0.44 ug/L 0.26 ug/L 0.29 ug/L	Qvalue 98 95 94 99

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5971.D RUN524.M Fri Feb 05 09:26:02 1999

Vial: 12

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V5971.D

Acq On : 4 Feb 99 6:10 pm

Sample : R-6171.3

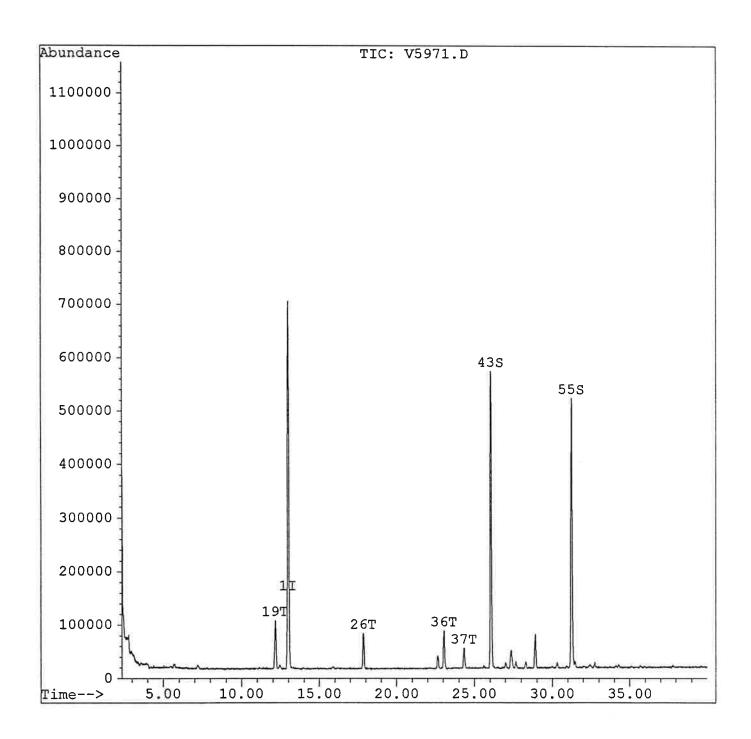
: Arco - RC-EE-64-0299

Misc

Quant Time: Feb 5 9:25 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V5972.D Vial: 13 Acq On : 4 Feb 99 6:57 pm Operator: vb

Inst : 5971 - In

Sample : R-6171.4 Misc : Arco - RC-EL-36-0299 Multiplr: 1.00

Quant Time: Feb 5 9:26 1999

Method Method : C:\HPCHEM\1\method \tag{\text{Non-in-obs}} \tag{\text{Non-in-obs}} \tag{\text{Title}} : 524.2 Purgable Organics : C:\HPCHEM\1\METHODS\RUN524.M Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.98	96	1970259	5.00 ug/L	0.00
System Monitoring Compounds				81	Recovery
43) 4-bromofluorobenzene	26.06	95	691063	4.82 ug/L	96.44%
55) 1,2-dichlorobenzene-d4	31.24	152	471666	5.00 ug/L	99.92%
Target Compounds					Qvalue
19) Benzene	12.15	78	183360	0.49 ug/L	96
26) Toluene	17.85	91	303069	0.82 ug/L	99
35) Ethylbenzene	22.66	91	127113	0.32 ug/L	95
36) m&p-xylenes	23.05	106	179066	0.63 ug/L	98
37) o-xylene	24.33	91	207837	0.74 ug/L	96

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5972.D RUN524.M Fri Feb 05 09:27:03 1999

Vial: 13

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V5972.D

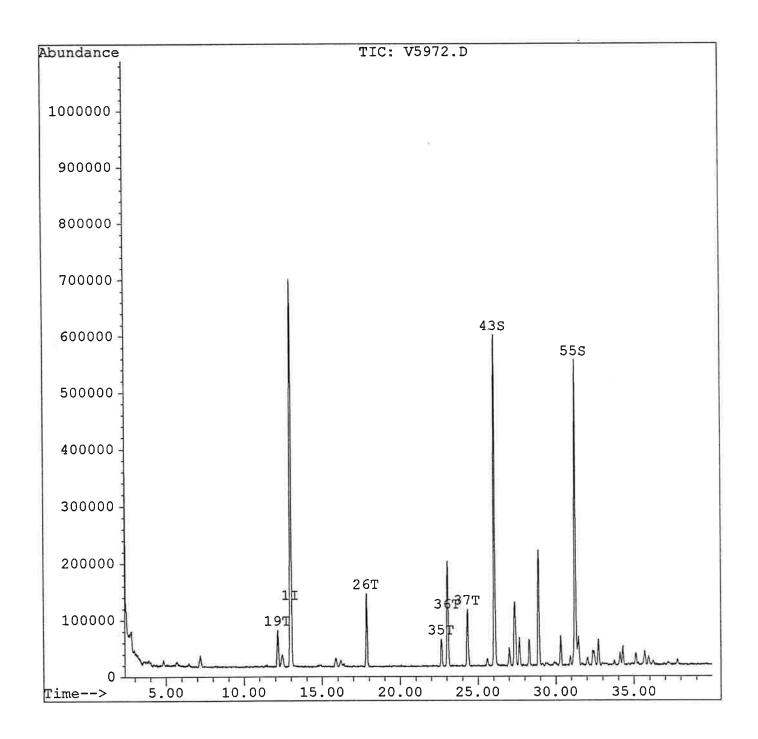
6:57 pm Acq On : 4 Feb 99

: R-6171.4 Sample

Misc : Arco - RC-EL-36-0299 Quant Time: Feb 5 9:26 1999

: C:\HPCHEM\1\METHODS\RUN524.M Method

: 524.2 Purgable Organics Title Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V5973.D Vial: 14 Operator: vb

Inst : 5971 - In

Acq On : 4 Feb 99 7:43 pm Sample : R-6171.5 Misc : Arco - RC-EL-18-0299 Multiplr: 1.00

Quant Time: Feb 5 9:28 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.99	96	1909289	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.05 31.24	95 152	669501 454757	%1 4.82 ug/L 4.97 ug/L	Recovery 96.41% 99.41%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	12.16 17.84 22.65 23.04 24.32	78 91 91 106 91	134071 279572 105827 153910 166791	0.37 ug/L 0.78 ug/L 0.27 ug/L 0.56 ug/L 0.61 ug/L	98 88

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5973.D RUN524.M Fri Feb 05 09:28:20 1999

Vial: 14

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V5973.D

Acq On : 4 Feb 99 7:43 pm

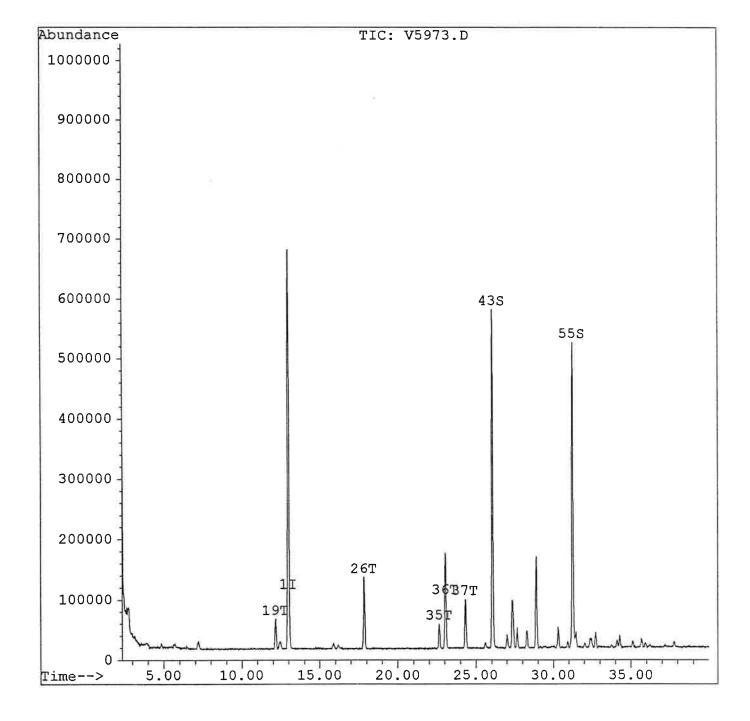
Sample : R-6171.5 Misc : Arco - RC-

: Arco - RC-EL-18-0299

Quant Time: Feb 5 9:28 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update: Thu Feb 04 16:27:27 1999
Response via: Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V5974.D

Vial: 15 Operator: vb

Inst : 5971 - In

Acq On : 4 Feb 99 8:29 pm Sample : R-6171.6 Misc : Arco - RC-EC-00-0299

Multiplr: 1.00

Quant Time: Feb 5 9:28 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.98	96	1928540	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.05 31.24	95 152	676570 468722	%I 4.82 ug/L 5.07 ug/L	
Target Compounds 19) Benzene 26) Toluene	12.16 17.84	78 91	214728 84547	0.58 ug/L 0.23 ug/L	Qvalue 99 93

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5974.D RUN524.M Fri Feb 05 09:29:09 1999

Vial: 15

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V5974.D

Acq On : 4 Feb 99 8:29 pm

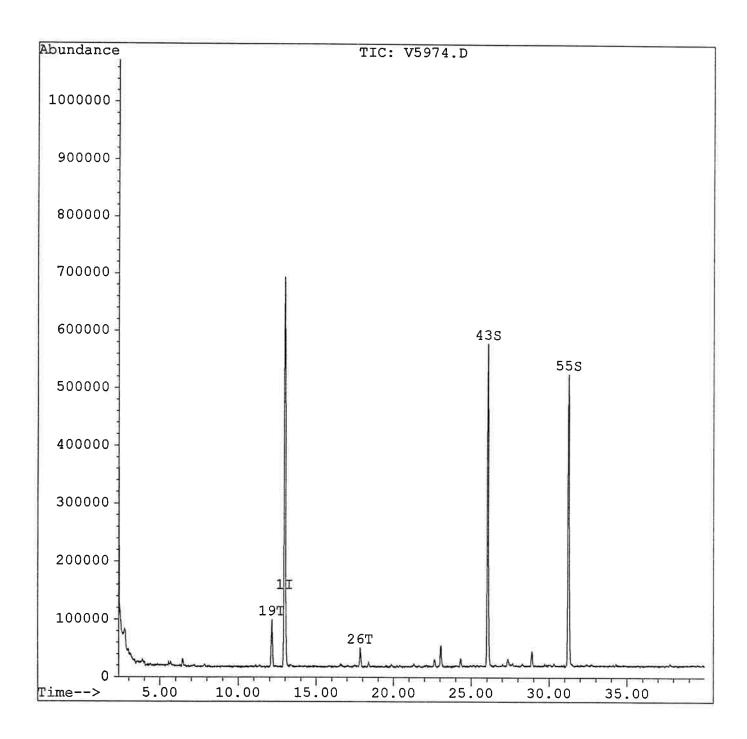
Sample : R-6171.6 Misc : Arco - RO

: Arco - RC-EC-00-0299

Quant Time: Feb 5 9:28 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



Data File : C:\HPCHEM\1\DATA\V5975.D

Vial: 16

Acq On : 4 Feb 99 9:16 pm Sample : R-6171.7 Misc : Arco - RC-EC-00-0299A

Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Quant Time: Feb 5 9:16 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.98	96	1867751	5.00 ug/L	0.00
System Monitoring Compounds				%R	ecovery
43) 4-bromofluorobenzene	26.06	95	653313	4.81 ug/L	96.17%
55) 1,2-dichlorobenzene-d4	31.23	152	437372	4.89 ug/L	97.74%
Target Compounds				#2	Qvalue
19) Benzene	12.16	78	229422	0.64  ug/L	100
26) Toluene	17.84	91	189861	0.54 ug/L	98
36) m&p-xylenes	23.04	106	80473	0.30 ug/L	92
37) o-xylene	24.33	91	81703	0.31 ug/L	99

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5975.D RUN524.M Fri Feb 05 09:30:05 1999

Vial: 16

: 5971 - In

Operator: vb

Multiplr: 1.00

Inst

Data File : C:\HPCHEM\1\DATA\V5975.D

Acq On : 4 Feb 99 9:16 pm

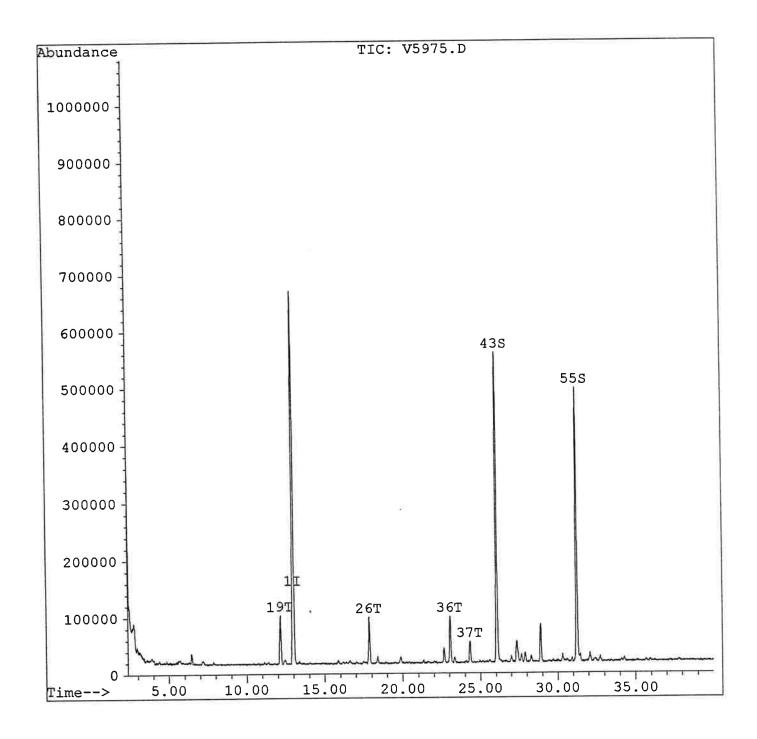
Sample : R-6171.7

Misc : Arco - RC-EC-00-0299A

Quant Time: Feb 5 9:16 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



Vial: 1 Data File : C:\HPCHEM\1\DATA\V5976.D Operator: vb Acq On : 4 Feb 99 10:02 pm

Inst : 5971 - In Sample : R-6171.8 Misc : Arco - RC-EC-34-0299

Multiplr: 1.00

Quant Time: Feb 5 9:15 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.98	96	1937935	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.05 31.24	95 152	676154 462193	%R 4.80 ug/L 4.98 ug/L	ecovery 95.93% 99.54%
Target Compounds 19) Benzene 26) Toluene 35) Ethylbenzene 36) m&p-xylenes 37) o-xylene	12.15 17.85 22.65 23.04 24.32	78 91 91 106 91	236137 220589 82917 100977 106963	0.64 ug/L 0.61 ug/L 0.21 ug/L 0.36 ug/L 0.39 ug/L	Qvalue 98 94 92 95 97

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5976.D RUN524.M Fri Feb 05 09:30:38 1999

Vial: 1 Operator: vb

Multiplr: 1.00

: 5971 - In

Inst

Data File : C:\HPCHEM\1\DATA\V5976.D

Acq On : 4 Feb 99 10:02 pm

: R-6171.8

: Arco - RC-EC-34-0299

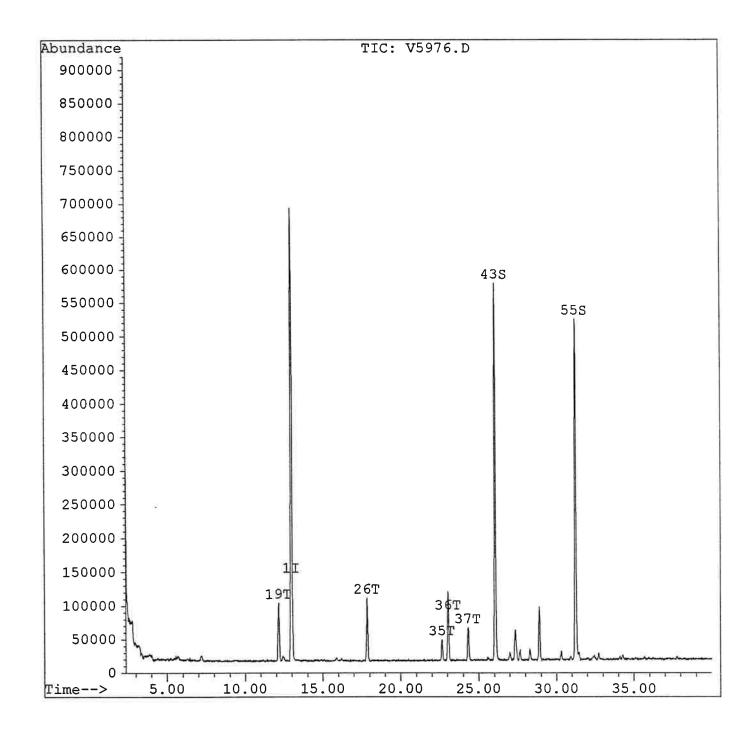
Quant Time: Feb 5 9:15 1999

Sample

Misc

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



Data File: C:\HPCHEM\1\DATA\V5977.D

Vial: 2 Operator: vb

Inst : 5971 - In

Acq On : 4 Feb 99 10:48 pm Sample : R-6171.9 Misc : Arco - Field Blank

Multiplr: 1.00

Quant Time: Feb 5 9:31 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics

Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	12.98	96	1820261	5.00 ug/L	0.00
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.05 31.23	95 152	647171 440825		97.76% 101.08%
Target Compounds 26) Toluene 36) m&p-xylenes	17.84 23.03	91 106	215081 52842	0.63 ug/L 0.20 ug/L	Qvalue 93 87

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5977.D RUN524.M Fri Feb 05 09:31:37 1999

Vial: 2

Multiplr: 1.00

Inst : 5971 - In

Operator: vb

Data File : C:\HPCHEM\1\DATA\V5977.D

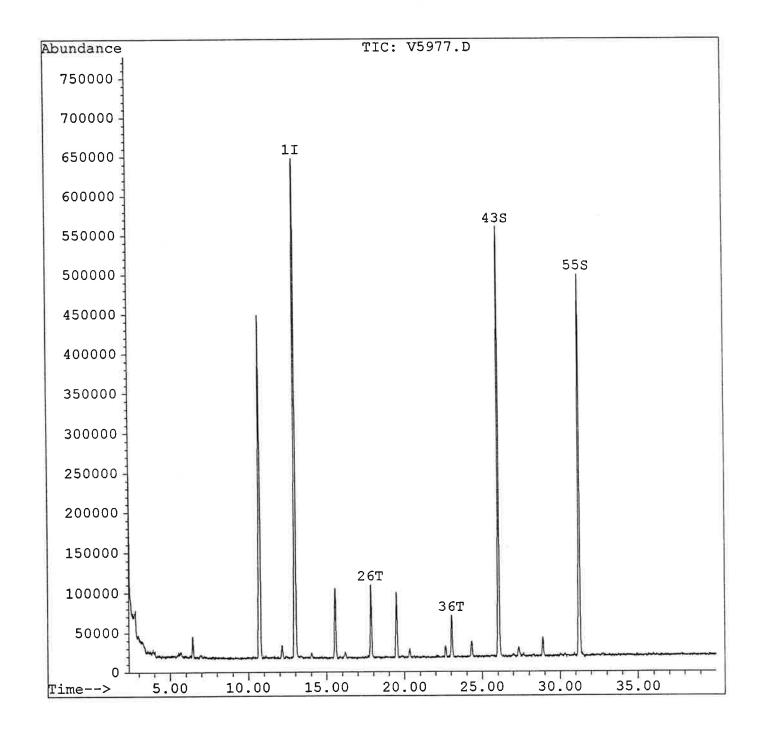
Acq On : 4 Feb 99 10:48 pm

Sample : R-6171.9 Misc : Arco - Fi

Misc : Arco - Field Blank Quant Time: Feb 5 9:31 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



### WATER VOLATILE SYSTEM MONITORING COMPOUND RECOVERY

Customer:	Arco

Γ		SMC1	SMC2		OTHER	TOT
L	SAMPLE NO.	#	#	#	#	OUT
01[	VBLK01	90	86			
02[	R-6171.1	98	95			
03[	R-6171.2	97	100			
04[	R-6171.3	96	99			
05[	R-6171.4	96	100			
06[	R-6171.5	96	99			
07[	R-6171.6	96	101			
08[	R-6171.7	96	98			
09[	R-6171.8	96	100			
10[	R-6171.9	98	101			
11[						
12[						
13[						
14						
15[						
16[						
17						
18						
19						
20[						
21						
22						
23[						
24[						
25[						
26						
27[						
28[						
29[						
30[						

QC LIMITS

(75-115)

(75-115)

SMC1 = 4-Bromofluorobenzene SMC2 = 1,2-dichlorobenzene-d4

# Column to be used to flag recovery values

- \* Values outside of contract required QC limits
- D System Monitoring Compound diluted out

FORM II VOA-1

### RELIANCE LABORATORIES, INC. WATER VOLATILE MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Matrix Spike - Sample No.: , R-6171.1

	SPIKE	SAMPLE	MS	MS	QC.
	ADDED	CONCENTRATION	CONCENTRATION	%	LIMITS
COMPOUND	(ug/Kg)	(ug/Kg)	(ug/Kg)	REC #	
Benzene	3.00	0.00	3.15	105	(80-120)
Toluene	3.00	0.00	2.97	99	(80-120)
Ethylbenzene	3.00	0.00	3.06	102	(80-120)
m&p-xylenes	3.00	0.00	3.08	103	(80-120)
o-xylenes	3.00	0.00	2.98	99	(80-120)
Styrene	3.00	0.00	3.11	104	(80-120)

	SPIKE ADDED	ADDED CONCENTRATION		%	QC LIMITS RPD   REC.	
COMPOUND	(ug/Kg)	(ug/Kg)	REC #	RPD #	RPD	
Benzene	3.00	3.06	102	3	20	(80-120)
Toluene	3.00	3.04	101	2	20	(80-120)
Ethylbenzene	3.00	3.14	105	3	20	(80-120)
m&p-xylenes	3.00	3.14	105	2	20	(80-120)
o-xylenes	3.00	3.06	102	3	20	(80-120)
Styrene	3.00	3.19	106	3	20	(80-120)

# Column to be used to flag recovery and RPD values with an asterisk
\* Values outside of QC limits

Comments:	

### **VOLATILE METHOD BLANK SUMMARY**

					VBLK
	Customer:	Arco	_		
Lab File ID V59	968.D			Lab Sample II	D: BLANK
Date Analyzed:	2/4/99			Time Analyze	d: 1552
GC Column:	DB-624 ID:	0.53 (mm	١		
	=======================================	(,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,,	,		
Instrument ID:		-			
THI	S METHOD BLANK	APPLIES TO THE	FOLLOWING	SAMPLES, MS	AND MSD:
		LAB	LAB	TIME	1
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	
	R-6171.1	ER-12	V5969.D	1639	1
	R-6171.2	ER-25	V5970.D	1724	]
	R-6171.3	EC-64	V5971.D	1810	
	R-6171.4	EL-36	V5972.D	1857	]
	R-6171.5	EL-18	V5973.D	1943	]
	R-6171.6	EC-00	V5974.D	2029	
	R-6171.7	EC-00A	V5975.D	2116	
	R-6171.8	EC-34	V5976.D	2202	]
	R-6171.9	FB	V5977.D	2248	]
10					
11					
12					
13					
14					1
15					1
16					
17					ĺ
18					
19					
20					
21					
22					
23					
24					
25					
26					
27					
28					
29					
30					

COMMENTS:		

Page 1 of 1

FORM IV VOA

Data File : C:\HPCHEM\1\DATA\V5968.D

Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Vial: 9

Acq On : 4 Feb 99 3:52 pm Sample : blank Misc : blank

Quant Time: Feb 5 9:19 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M
Title : 524.2 Purgable Organics Last Update : Thu Feb 04 16:27:27 1999 Response via : Multiple Level Calibration

Internal Standards	R.T.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene	13.00	96	1972619	5.00 ug/I	0.01
System Monitoring Compounds 43) 4-bromofluorobenzene 55) 1,2-dichlorobenzene-d4	26.06 31.26	95 152	644818 404960	4.49 ug/l 4.28 ug/l	
Target Compounds					Qvalue

<sup>(#) =</sup> qualifier out of range (m) = manual integration V5968.D RUN524.M Fri Feb 05 09:19:24 1999

Vial: 9

Multiplr: 1.00

: 5971 - In

Operator: vb

Inst

Data File : C:\HPCHEM\1\DATA\V5968.D

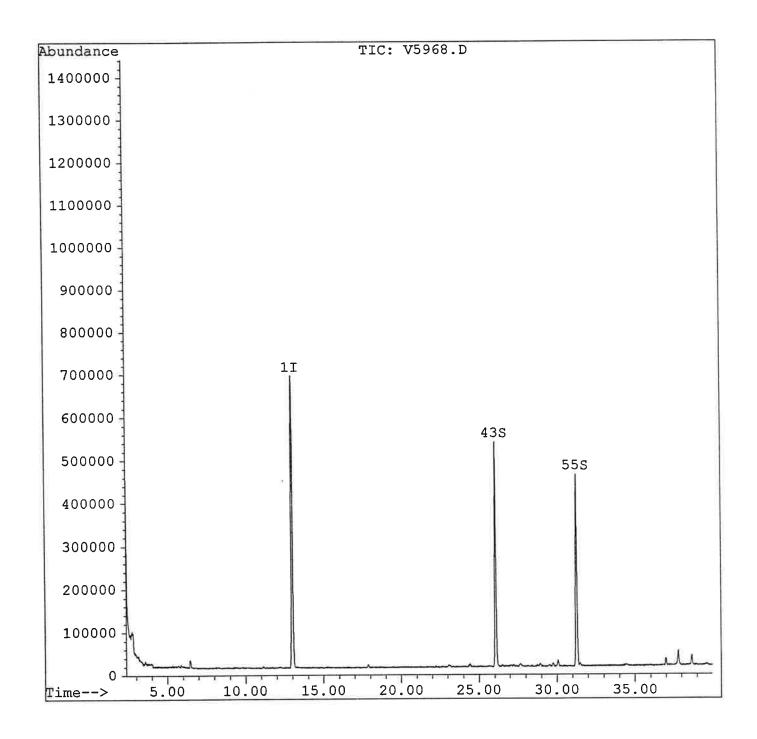
Acq On : 4 Feb 99 3:52 pm

Sample : blank
Misc : blank

Quant Time: Feb 5 9:19 1999

Method : C:\HPCHEM\1\METHODS\RUN524.M

Title : 524.2 Purgable Organics
Last Update : Thu Feb 04 16:27:27 1999
Response via : Multiple Level Calibration



### VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK BROMOFLUOROBENZENE (BFB)

Customer	:	Arco		

Lab File ID: V5964.D BFB Injection Date: 2/4/99

Instrument ID: HP5971A BFB Injection Time: 1134

GC Column: \_\_DB-624 ID: 0.53 (mm)

	%RELATIVE		
m/e	ION ABUNDANCE CRITERIA	ABUNDANCE	
50	8.0 - 40.0% of mass 95	19.5	
75	30.0 - 66.0% of mass 95	42.2	
95	Base peak, 100% relative abundance	100.0	
96	5.0 - 9.0% of mass 95	6.2	
173	Less than 2.0% of mass 174	0.0 (	0.0)1
174	50.0 - 120.0% of mass 95	86.6	
175	4.0 - 9.0% of mass 174	6.4 (	7.4)1
176	93.0 - 101.0% of mass 174	84.8 (	7.9 )1
177	5.0 - 9.0% of mass 176	5.5 (	6.4)2

1-Value is % mass 174 2-Value is % mass 176

This check applies to the following SAMPLES, MS, MSD, BLANKS and STANDARDS:

	LAB	LAB	DATE	TIME
SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED	ANALYZED
01 VSTD010	ICC001	V5965.D	2/4/99	1217
02 VSTD020	ICC002	V5966.D	2/4/99	1304
03 VSTD050	ICC005	V5967.D	2/4/99	1505
VBLK01	BLANK	V5968.D	2/4/99	1552
05 R-6171.1	ER-12	V5969.D	2/4/99	1639
06 R-6171.2	ER-25	V5970.D	2/4/99	1724
7 R-6171.3	EC-64	V5971.D	2/4/99	1810
08 R-6171.4	EL-36	V5972.D	2/4/99	1857
9 R-6171.5	EL-18	V5973.D	2/4/99	1943
IO R-6171.6	EC-00	V5974.D	2/4/99	2029
I1 R-6171.7	EC-00A	V5975.D	2/4/99	2116
I2 R-6171.8	EC-34	V5976.D	2/4/99	2202
I3 R-6171.9	FB	V5977.D	2/4/99	2248
14				
15				
16				
17				
18				
19				
20				
21				
22				

Page 1 of 1

FORM V VOA

Data File : C:\HPCHEM\1\DATA\V5964.D

Acq On : 4 Feb 99 11:34 am

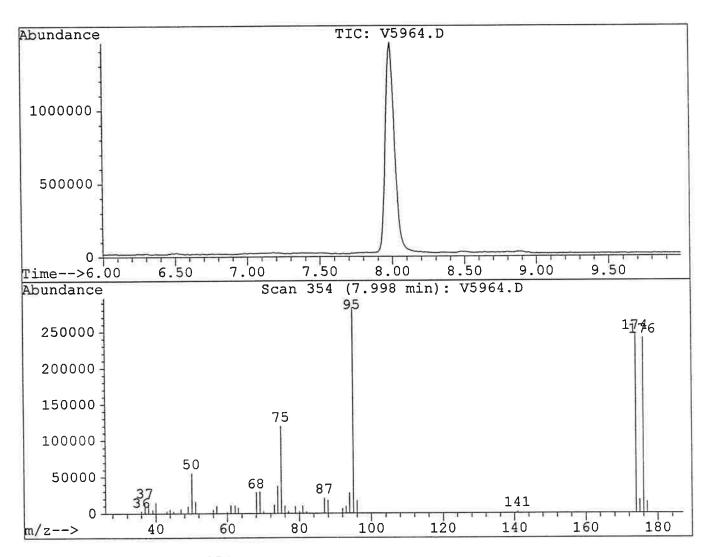
Sample : bfb Misc : bfb Vial: 1
Operator: vb

Inst : 5971 - In

Multiplr: 1.00

Method : C:\HPCHEM\1\METHODS\ENVDEF.M

Title :



Peak Apex is scan: 354

1	Target Mass	 	Rel. to Mass	1	Lower Limit%	1	Upper Limit%	1	Rel. Abn%	l l	Raw Abn	1	Result Pass/Fail	
1	50	ı	95	1	15	1	40	1	19.5	1	55304	1	PASS	1
i	75	i	95	ĺ	30	Î	60	1	42.2	ı	119584	1	PASS	1
î	95	ì	95	î	100	1	100	I	100.0	1	283136	1	PASS	1
ì	96	ï	95	Ĺ	5	1	9	1	6.2	1	17432	1	PASS	1
î	173	ï	174	i	0	1	2	1	0.0	Ĩ	0	1	PASS	1
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î	175	i	174	i	5	1	9	1	7.4	Ì	18120	1	PASS	1
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### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Custo	mer Arco		
Instrument ID: HP5971A	Calibration Date(s):	2/4/99	2/4/99
	Calibration Times:	1217	1505
GC Column: DB-624	ID: <u>0.53</u> (mm)		

Lab File ID: RRF05 = V5967.D	RRF01 =	V5965.D		RRF02 =	V5966.D		
						RRF	% RSE
COMPOUND	RRF01	RRF02	RRF05				
Benzene	0.986	0.944	0.944			0.958	2.5
Toluene	1.030	0.879	0.902			0.937	8.7
Ethylbenzene	1.070	0.954	1.003			1.009	5.8
m&p-xylenes	0.759	0.687	0.728			0.725	5.0
o-xylene	0.763	0.662	0.705			0.710	7.1
Styrene	0.547	0.480	0.504			0.510	6.7
4-bromofluorobenzene	0.370	0.367	0.354			0.364	2.3
1,2-dichlorobenzene-d4	0.239	0.247	0.232			0.239	3.1

Page 1 of 1

FORM VI VOA

### VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Customer : Arco

Lab File ID (Standard): V5967.D

Date Analyzed: 2/4/99

Instrument ID: HP5971A

Time Analyzed: \_\_\_1505

GC Column: DB-624

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ID: 0.53 (mm)

	IS1									
	AREA #	RT #	AREA	#	RT	#	AREA	#	RT	#
12 HOUR STD	2127714	13.00						-		
UPPER LIMIT	4255428	13.50								
LOWER LIMIT	1063857	12.50								
SAMPLE										
NO.										
01 VBLK01	1972619	13.00								
02 R-6171.1	1967274	12.99								
03 R-6171.2	1877487	12.99								
04 R-6171.3	1941054	12.99								
05 R-6171.4	1970259	12.98								
06 R-6171.5	1909289	12.99								
07 R-6171.6	1928540	12.98								
08 R-6171.7	1867751	12.98								
09 R-6171.8	1937935	12.98								
10 R-6171.9	1820261	12.98								
11										
12										
13										
14										
15										
16										
17										
18										
19										
20										
21										
22										

IS1 = Fluorobenzene

AREA UPPER LIMIT = +100% of internal standard area
AREA LOWER LIMIT = -50% of internal standard area
RT UPPER LIMIT = +0.50 minutes of internal standard RT
RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk

\* Values outside of QC limits.

Page 1 of 1

FORM VIII VOA



### DEPARTMENT OF ENVIRONMENTAL PROTECTION

Certifies That
Reliance Laboratories, Inc.
3090 Wood Bridge Avenue
Edison, NJ 08837



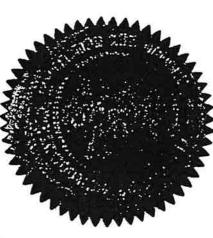
having duly met the requirements of the

Regulations Governing Laboratory Certification And Standards Of Performance NJ.A.C. 7:18 et. seq.

is hereby approved as a

## State Certified Water Laboratory

To perform the analyses as indicated on the Annual Certified Parameter List which must accompany this certificate to be valid



# 12687 PERMANENT CERTIFICATION NUMBER

anthe ! Day with

N.J.A.C. 7:18-2.11(d) and agreed to by the Laboratory Manager on filing the application This certification is subject to unannounced laboratory inspections as specified by

TO BE CGNSPICUOUSLY DISPLAYED AT THE LABORATORY WITH THE ANNUAL CERTIFIED PARAMETER LIST.

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